

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE082815\
 Data File : BE090656.D
 Acq On : 28 Aug 2015 11:51
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001

Manual Integrations
 APPROVED

MMDadoda
 8/31/2015 1:46:38 PM

Quant Time: Aug 31 04:57:43 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE081415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 14 16:52:42 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	22013	5.00	ng	0.00
6) Naphthalene-d8	10.33	136	100368	5.00	ng	0.00
12) Acenaphthene-d10	14.19	164	58832	5.00	ng	0.00
18) Phenanthrene-d10	16.94	188	153497	5.00	ng	0.00
25) Chrysene-d12	21.09	240	164256	5.00	ng	0.00
32) Perylene-d12	23.39	264	136988	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.21	112	3934m	0.62	ng	0.00
5) Phenol-d6	6.78	99	5870	0.63	ng	0.00
7) Nitrobenzene-d5	8.73	82	7085	0.89	ng	0.00
13) 2,4,6-Tribromophenol	15.71	330	1425	0.63	ng	0.00
14) 2-Fluorobiphenyl	12.82	172	15485	0.88	ng	0.00
27) Terphenyl-d14	19.56	244	25888	0.86	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.19	88	3332	1.28	ng	95
3) n-Nitrosodimethylamine	3.48	42	3751	0.99	ng	# 86
8) Nitrobenzene	8.77	77	6984	0.84	ng	97
9) Naphthalene	10.38	128	22557	0.99	ng	99
10) Hexachlorobutadiene	10.67	225	3630	1.08	ng	99
11) 2-Methylnaphthalene	12.01	142	13236	0.89	ng	97
15) Acenaphthylene	13.91	152	94781	0.88	ng	100
16) Acenaphthene	14.25	154	15871	0.98	ng	84
17) Fluorene	15.26	166	20738	1.05	ng	99
19) 4-Bromophenyl-phenylether	16.14	248	5621	0.90	ng	83
20) Hexachlorobenzene	16.25	284	28003	1.01	ng	97
21) Pentachlorophenol	16.61	266	1220	0.65	ng	95
22) Phenanthrene	16.98	178	38599	0.96	ng	100
23) Anthracene	17.07	178	32946	0.89	ng	100
24) Fluoranthene	18.98	202	39175	0.91	ng	99
26) Pyrene	19.34	202	41167	0.88	ng	100
28) Benzo(a)anthracene	21.07	228	38566	0.89	ng	99
29) Chrysene	21.13	228	43824	0.99	ng	99
30) Bis(2-ethylhexyl)phthalate	21.02	149	9449	0.50	ng	99
31) Indeno(1,2,3-cd)pyrene	25.78	276	33034	0.79	ng	# 92
33) Benzo(b)fluoranthene	22.70	252	30400	0.94	ng	99
34) Benzo(k)fluoranthene	22.74	252	36682	1.00	ng	99
35) Benzo(a)pyrene	23.29	252	26955	0.89	ng	# 96
36) Dibenzo(a,h)anthracene	25.81	278	24131	0.82	ng	97
37) Benzo(g,h,i)perylene	26.51	276	27416	0.86	ng	95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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